chain nodes : 7 8 9 10 11 12 13 15 17 18 29 30 31 32 33 34 ring nodes : 1 2 3 4 5 6 14 16 19 20 21 22 23 24 25 26 27 28 chain bonds : 2-7 5-13 7-8 8-9 8-10 8-11 10-34 11-12 13-14 13-15 15-16 15-17 17-18 26-29 29-30 30-31 31-32 31-33 ring bonds : 1-2 1-6 2-3 3-4 4-5 5-6 14-24 14-28 16-19 16-23 19-20 20-21 21-22 22-23 24-25 25-26 26-27 27-28 exact/norm bonds : 2-7 7-8 31-32 31-33 exact bonds : 5-13 10-34 11-12 13-14 13-15 15-16 15-17 17-18 26-29 29-30 30-31 normalized bonds : 1-2 1-6 2-3 3-4 4-5 5-6 8-9 8-10 8-11 14-24 14-28 16-19 16-23 19-20 20-21 21-22 22-23 24-25 25-26 26-27 27-28

## Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:Atom 16:CLASS 17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:Atom 30:Atom 31:Atom 32:Atom 34:Atom

## L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

=> d L2

L2 HAS NO ANSWERS

L2 ST

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

Structure attributes must be viewed using STN Express query preparation.

=> d L1

L1 HAS NO ANSWERS

L1 SCR 963 AND 1006 AND 2076

=> s L2 full

FULL SEARCH INITIATED 14:48:20 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 64 TO ITERATE

100.0% PROCESSED 64 ITERATIONS

2 ANSWERS

SEARCH TIME: 00.00.06

L4 2 SEA SSS FUL L2

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 172.55 172.76

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 14:48:30 ON 20 AUG 2007 .
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=> s L4

L5 1 L4

=> d L5 bib abs hitstr

- L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN
- AN 2005:324121 CAPLUS
- DN 142:392179
- TI Preparation of acyloxydiphenylbutenylcinnamates as estrogen receptor modulator prodrugs.
- IN <u>Eaddy</u>, John Fred, III; Heyer, Dennis; Katamreddy, Subba Reddy; Martin, Michael Tolar; McClure, Michael Scott; Randhawa, Amarjit Sab; Samano, Vicente; Ray, John Albert

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PA
     Smithkline Beecham Corporation, USA
SO
     PCT Int. Appl., 78 pp.
     CODEN: PIXXD2
DT
     Patent
LA
     English
FAN.CNT 1
     PATENT NO.
                          KIND
                                 DATE
                                              APPLICATION NO.
                                                                      DATE
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ΡI
     WO 2005033056
                          A2
                                  20050414
                                              WO 2004-US32918
                                                                       20041004
     WO 2005033056
                          A3
                                  20050623
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     US 2007111971
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                                  20070517
PRAI US 2003-509678P
                           Р
                                  20031008
                           Р
     US 2003-514692P
                                  20031027
     WO 2004-US32918
                           W
                                  20041004
OS
     CASREACT 142:392179; MARPAT 142:392179
GI
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CN

AB Title compds. (I; R1 = ACO, PO3H2; A = alkyl, aryl, heteroaryl, cycloalkyl, aminoalkyl, alkoxy, alkoxyalkyl, haloalkyl, heterocyclylalkyl), were prepared Thus, I (R1 = H) (preparation given) and Et3N in THF at 5° were treated with propionyl chloride in THF followed by stirring for 1 h to give 64% I (R1 = EtCO). The latter orally in rats showed 86.6% bioavailability, vs. 5.7% for I (R1 = H). IT 850005-29-7P RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (claimed compound; preparation of acyloxydiphenylbutenylcinnamates as estrogen receptor modulator prodrugs) RN 850005-29-7 CAPLUS

2-Propenoic acid, 3-[4-[(1Z)-2-phenyl-1-[4-(phosphonooxy)phenyl]-1-

butenyl]phenyl]-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

IT 850005-46-8P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of acyloxydiphenylbutenylcinnamates as estrogen receptor modulator prodrugs)

RN 850005-46-8 CAPLUS

CN 2-Propenoic acid, 3-[4-[(1Z)-2-phenyl-1-[4-(phosphonooxy)phenyl]-1-butenyl]phenyl]-, 1-(1,1-dimethylethyl) ester, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

=>

---Logging off of STN---

=>

Executing the logoff script...

chain nodes : 7 9 10 11 23 24 25 26 27 ring nodes : 1 2 3 4 5 6 8 12 13 14 15 16 17 18 19 20 21 22 chain bonds : 1-7 4-25 7-8 7-9 9-10 9-12 10-11 20-23 23-24 25-26 26-27 ring bonds : 1-2 1-6 2-3 3-4 4-5 5-6 8-18 8-22 12-13 12-17 13-14 14-15 15-16 16-17 18-19 19-20 20-21 21-22 exact/norm bonds : 20-23 23-24 exact bonds : 1-7 4-25 7-8 7-9 9-10 9-12 10-11 25-26 26-27 normalized bonds : 1-2 1-6 2-3 3-4 4-5 5-6 8-18 8-22 12-13 12-17 13-14 14-15 15-16 16-17 18-19 19-20 20-21 21-22

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## L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

=> d L2

L2 HAS NO ANSWERS